KRS 10/5/2007

EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	2713	ab initio	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:21
L2	1973	interaction energy	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:22
L3	2	l2 with l1	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:22
L4	281	molecular fragmentation	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:22
L5	9	molecular cap	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:23
L6	1541	molecular mechanics	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:23
L7	19	l2 with l6	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:27

EAST Search History

L8	1	molecular interaction energy with calculation	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:29
L9	160	interaction energy with calculation	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 15:50
L10	26	l1 and l9	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 14:29
L12	136	hartree-fock	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 15:51
L13	11	l12 and l2	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT; IBM_TDB	ADJ	ON	2007/10/05 15:52

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     (FILE 'HOME' ENTERED AT 14:03:33 ON 05 OCT 2007)

     FILE 'MEDLINE, CAPLUS, EMBASE, BIOSIS, PASCAL, SCISEARCH' ENTERED AT
     14:04:14 ON 05 OCT 2007
L1             0 S QUATUM MECHANICAL INTERMOLEULAR INTERACTION ENERGY
L2             0 S QUANTUM MECHANICAL INTERMOLEULAR INTERACTION ENERGY
L3            25 S QUANTUM MECHANICAL INTERACTION ENERGY
L4            14 DUP REM L3 (11 DUPLICATES REMOVED)
L5          8342 S (AB INITIO) AND (QUANTUM MECHANICAL)
L6        279104 S FRAGMENTATION
L7            57 S L5 AND L6
L8            59 S MOLECULAR CAP
L9             0 S L7 AND L8
L10            5 S L5 AND L8
L11            3 DUP REM L10 (2 DUPLICATES REMOVED)
L12            1 S L5 AND (GHOST ATOM)
L13         2872 S (AB INITIO) AND (QUANTUM MECHANIC)
L14           12 S L6 AND L13
L15            8 DUP REM L14 (4 DUPLICATES REMOVED)
L16          136 S AB INITIO ENERGY CALCULATION
L17           44 S L16 AND INTERACTION
L18           15 DUP REM L17 (29 DUPLICATES REMOVED)

     FILE 'STNGUIDE' ENTERED AT 14:18:29 ON 05 OCT 2007
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